

CLAIMS

We claim:

1. A computer implemented method to search a heterogenous compound database for molecules which are likely to have the same biological activity as a known query molecule comprising the following steps:
 - a) defining fragments of a query molecule and a database molecules according to a defined set of rules;
 - b) generating shape descriptors for the query molecule and database molecule fragments; and
 - c) using the shape descriptors, identifying the database molecule which has a shape similar to the query molecule.
- 10 2. A computer implemented method to search a heterogenous compound database for molecules which are likely to have the same biological activity as a known query molecule comprising the following steps:
 - a) fragmenting a query molecule according to a defined set of rules;
 - b) topomerically aligning the query molecule fragments to generate a topomeric conformation;
 - c) generating the interaction energies between a probe and the atoms in the topomerically aligned query fragments at all intersection points in a three dimensional grid surrounding the aligned query fragments;
 - 15 d) fragmenting a database compound according to a defined set of rules;
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- e) topomerically aligning the database molecule fragments to generate a topomeric conformation;
 - f) generating the interaction energies between a probe and the atoms in the topomerically aligned database fragments at all intersection points in a three dimensional grid surrounding the aligned query fragments;
 - g) determining the similarity between query and database fragments by the root sum square differences in the field values; and
 - h) identify the molecule in the database most similar to the query molecule as that molecule having the smallest field value difference in its fragments.

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